

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022522\
 Data File : BM034090.D
 Acq On : 25 Feb 2022 23:14
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 26 04:02:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 26 03:45:21 2022
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	59	0.00
2	1,4-Dioxane	40.000	38.066	4.8	57	0.00
3	Pyridine	40.000	37.016	7.5	54	0.00
4	n-Nitrosodimethylamine	40.000	39.394	1.5	58	-0.01
5 S	2-Fluorophenol	80.000	78.206	2.2	57	0.00
6	Aniline	40.000	36.851	7.9	53	0.00
7 S	Phenol-d6	80.000	75.003	6.2	54	0.00
8	2-Chlorophenol	40.000	39.536	1.2	56	0.00
9	Benzaldehyde	40.000	34.536	13.7	56	0.00
10 C	Phenol	40.000	36.710	8.2	53	0.00
11	bis(2-Chloroethyl)ether	40.000	35.739	10.7	52	0.00
12	1,3-Dichlorobenzene	40.000	40.032	-0.1	58	0.00
13 C	1,4-Dichlorobenzene	40.000	40.135	-0.3	59	0.00
14	1,2-Dichlorobenzene	40.000	39.779	0.6	58	0.00
15	Benzyl Alcohol	40.000	38.708	3.2	56	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	30.767	23.1	45	0.00
17	2-Methylphenol	40.000	38.110	4.7	55	0.00
18	Hexachloroethane	40.000	39.593	1.0	58	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	36.979	7.6	53	0.00
20	3+4-Methylphenols	40.000	37.836	5.4	54	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	56	0.00
22	Acetophenone	40.000	38.895	2.8	55	0.00
23 S	Nitrobenzene-d5	80.000	78.597	1.8	55	0.00
24	Nitrobenzene	40.000	38.340	4.1	54	0.00
25	Isophorone	40.000	38.167	4.6	53	0.00
26 C	2-Nitrophenol	40.000	44.163	-10.4	60	0.00
27	2,4-Dimethylphenol	40.000	38.193	4.5	54	0.00
28	bis(2-Chloroethoxy)methane	40.000	36.333	9.2	52	0.00
29 C	2,4-Dichlorophenol	40.000	41.807	-4.5	58	0.00
30	1,2,4-Trichlorobenzene	40.000	41.576	-3.9	60	0.00
31	Naphthalene	40.000	38.893	2.8	55	0.00
32	Benzoic acid	40.000	7.098	82.3#	10	-0.07
33	4-Chloroaniline	40.000	39.517	1.2	56	0.00
34 C	Hexachlorobutadiene	40.000	43.330	-8.3	62	0.00
35	Caprolactam	40.000	59.898	-49.7#	85	0.00
36 C	4-Chloro-3-methylphenol	40.000	41.137	-2.8	58	0.00
37	2-Methylnaphthalene	40.000	40.087	-0.2	57	0.00
38	1-Methylnaphthalene	40.000	40.273	-0.7	57	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	60	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	40.796	-2.0	61	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.253	-3.1	60	0.00
42 S	2,4,6-Tribromophenol	80.000	95.272	-19.1	75	0.00
43 C	2,4,6-Trichlorophenol	40.000	45.412	-13.5	68	0.00
44	2,4,5-Trichlorophenol	40.000	47.041	-17.6	71	0.00
45 S	2-Fluorobiphenyl	80.000	79.026	1.2	59	0.00
46	1,1'-Biphenyl	40.000	38.506	3.7	58	0.00
47	2-Chloronaphthalene	40.000	38.853	2.9	58	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022522\
 Data File : BM034090.D
 Acq On : 25 Feb 2022 23:14
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 26 04:02:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 26 03:45:21 2022
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	40.963	-2.4	58	0.00
49	Acenaphthylene	40.000	39.578	1.1	59	0.00
50	Dimethylphthalate	40.000	41.996	-5.0	64	0.00
51	2,6-Dinitrotoluene	40.000	43.233	-8.1	64	0.00
52 C	Acenaphthene	40.000	37.204	7.0	56	0.00
53	3-Nitroaniline	40.000	43.082	-7.7	63	0.00
54 P	2,4-Dinitrophenol	40.000	9.195	77.0#	9	0.00
55	Dibenzofuran	40.000	40.165	-0.4	61	0.00
56 P	4-Nitrophenol	40.000	35.822	10.4	52	0.00
57	2,4-Dinitrotoluene	40.000	45.312	-13.3	66	0.00
58	Fluorene	40.000	41.179	-2.9	62	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	43.355	-8.4	65	0.00
60	Diethylphthalate	40.000	42.812	-7.0	65	0.00
61	4-Chlorophenyl-phenylether	40.000	42.027	-5.1	64	0.00
62	4-Nitroaniline	40.000	44.553	-11.4	64	0.00
63	Azobenzene	40.000	38.225	4.4	58	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	66	0.00
65	4,6-Dinitro-2-methylphenol	40.000	8.893	77.8#	14	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.615	3.5	63	0.00
67	4-Bromophenyl-phenylether	40.000	41.400	-3.5	72	0.00
68	Hexachlorobenzene	40.000	41.248	-3.1	70	0.00
69	Atrazine	40.000	44.654	-11.6	71	0.00
70 C	Pentachlorophenol	40.000	29.872	25.3#	47	0.00
71	Phenanthrene	40.000	39.622	0.9	66	0.00
72	Anthracene	40.000	40.416	-1.0	66	0.00
73	Carbazole	40.000	40.789	-2.0	67	0.00
74	Di-n-butylphthalate	40.000	42.132	-5.3	67	0.00
75 C	Fluoranthene	40.000	42.455	-6.1	69	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	72	0.00
77	Benzidine	40.000	46.378	-15.9	68	0.00
78	Pyrene	40.000	37.491	6.3	70	0.00
79 S	Terphenyl-d14	80.000	79.505	0.6	73	0.00
80	Butylbenzylphthalate	40.000	39.673	0.8	68	0.00
81	Benzo(a)anthracene	40.000	39.971	0.1	72	0.00
82	3,3'-Dichlorobenzidine	40.000	43.545	-8.9	75	0.00
83	Chrysene	40.000	40.298	-0.7	72	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	39.845	0.4	67	0.00
85 c	Di-n-octyl phthalate	40.000	40.662	-1.7	68	0.00
86 I	Perylene-d12	20.000	20.000	0.0	75	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	41.510	-3.8	77	0.00
88	Benzo(b)fluoranthene	40.000	39.905	0.2	75	0.00
89	Benzo(k)fluoranthene	40.000	39.918	0.2	74	0.00
90 C	Benzo(a)pyrene	40.000	40.362	-0.9	75	0.00
91	Dibenzo(a,h)anthracene	40.000	41.284	-3.2	77	-0.01
92	Benzo(g,h,i)perylene	40.000	41.561	-3.9	77	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022522\
Data File : BM034090.D
Acq On : 25 Feb 2022 23:14
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Feb 26 04:02:59 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Feb 26 03:45:21 2022
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

Compound	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

(#) = Out of Range

SPCC's out = 0 CCC's out = 1